Searches for User lwells (Count = 3529)

Queries 3480 through 3529.

S #
Updt
Database
Query
Time
Comment

S3529
U
USPT
6416760.pn.
2002-11-08
07:54:12

S3528
U
USPT
irritant side effect
2002-11-07
16:13:54

S3527
U
USPT,JPAB,EPAB,DWPI
((424/401)!.CCLS.) and (psoriasis same (alpha keto acid or pyruvic or keto octanoic or beta
keto acid or acetoacetic or retinoid or retinoic or retinol or anthralin or anthranoid or
dioxyanthranol or peroxide or benzoyl peroxide or minoxidil or lithium salt or antimetabolite or
vitamin d or depigmentation or depigmenting or hydroquinone))
2002-11-07
12:50:50

S3526

U

USPT,JPAB,EPAB,DWPI

psoriasis same (alpha keto acid or pyruvic or keto octanoic or beta keto acid or acetoacetic or
retinoid or retinoic or retinol or anthralin or anthranoid or dioxyanthranol or peroxide or benzoyl
peroxide or minoxidil or lithium salt or antimetabolite or vitamin d or depigmentation or
depigmenting or hydroquinone)

2002-11-07
12:50:36

S3525

U

USPT,JPAB,EPAB,DWPI

(nsaid or antiinflammatory or anti inflammatory) near (alpha keto acid or pyruvic or keto
octanoic or beta keto acid or acetoacetic or retinoid or retinoic or retinol or anthralin or
anthranoid or dioxyanthranol or peroxide or benzoyl peroxide or minoxidil or lithium salt or
antimetabolite or vitamin d or depigmentation or depigmenting or hydroquinone)

2002-11-07
12:38:12

S3524

U

USPT,JPAB,EPAB,DWPI

((424/401)!.CCLS.) and ((nsaid or antiinflammatory or anti inflammatory) same (alpha keto
acid or pyruvic or keto octanoic or beta keto acid or acetoacetic or retinoid or retinoic or retinol
or anthralin or anthranoid or dioxyanthranol or peroxide or benzoyl peroxide or minoxidil or
lithium salt or antimetabolite or vitamin d or depigmentation or depigmenting or hydroquinone
))

2002-11-07
12:37:54

S3523

U

USPT,JPAB,EPAB,DWPI

(nsaid or antiinflammatory or anti inflammatory) same (alpha keto acid or pyruvic or keto
octanoic or beta keto acid or acetoacetic or retinoid or retinoic or retinol or anthralin or

anthranoid or dioxyanthranol or peroxide or benzoyl peroxide or minoxidil or lithium salt or
antimetabolite or vitamin d or depigmentation or depigmenting or hydroquinone)

2002-11-07
12:28:44

S3522
U
USPT,JPAB,EPAB,DWPI
sodium cromoglycate same (alpha keto acid or pyruvic or keto octanoic or beta keto acid or
acetoacetic or retinoid or retinoic or retinol or anthralin or anthranoid or dioxyanthranol or
peroxide or benzoyl peroxide or minoxidil or lithium salt or antimetabolite or vitamin d or
depigmentation or depigmenting or hydroquinone)

2002-11-07
12:26:40

S3521
U
USPT,JPAB,EPAB,DWPI
hair and (cosmetic and ((tnf near alpha near antagonist)not (tnf near alpha near antagonist)))

2002-11-07
12:21:59

S3520
U
USPT,JPAB,EPAB,DWPI
cosmetic and (((424/401)!.CCLS.) and ((tnf near alpha near antagonist)not (tnf near alpha near
antagonist)))

2002-11-07
12:21:41

S3519
U
USPT,JPAB,EPAB,DWPI
((424/401)!.CCLS.) and (sodium cromoglycate)

2002-11-07
12:21:26

S3518
U
USPT,JPAB,EPAB,DWPI
sodium cromoglycate

2002-11-07
12:21:15

S3517

U

USPT,JPAB,EPAB,DWPI

(((lisophyline or a802715 or sulphasalazine)and (pharmaceutical or cosmetic))not (((tnf near alpha near antagonist)not (tnf near alpha near antagonist))or ((interleukin 1 near antagonist)and (interleukin 1 near antagonist))or ((interleukin 1 near antagonist)same (interleukin 1 near antagonist))or (cosmetic and (alpha keto acid or pyruvic or keto octanoic or beta keto acid or acetoacetic or retinoid or retinoic or retinol or anthralin or anthranoid or dioxyanthranol or peroxide or benzoyl peroxide or minoxidil or lithium salt or antimetabolite or vitamin d or depigmentation or depigmenting or hydroquinone)))) not (((auranofin or skf-105809 or lactoferin)or (lisophyline or a802715 or sulphasalazine)or (interleukin 1 near antagonist)or (tnf near alpha near antagonist)) and (alpha keto acid or pyruvic or keto octanoic or beta keto acid or acetoacetic or retinoid or retinoic or retinol or anthralin or anthranoid or dioxyanthranol or peroxide or benzoyl peroxide or minoxidil or lithium salt or antimetabolite or vitamin d or depigmentation or depigmenting or hydroquinone))

2002-11-07 11:32:47

S3516

U

USPT,JPAB,EPAB,DWPI

((lisophyline or a802715 or sulphasalazine)and (pharmaceutical or cosmetic)) not (((tnf near alpha near antagonist)not (tnf near alpha near antagonist)) or ((interleukin 1 near antagonist)and (interleukin 1 near antagonist)) or ((interleukin 1 near antagonist)same (interleukin 1 near antagonist)) or (cosmetic and (alpha keto acid or pyruvic or keto octanoic or beta keto acid or acetoacetic or retinoid or retinoic or retinol or anthralin or anthranoid or dioxyanthranol or peroxide or benzoyl peroxide or minoxidil or lithium salt or antimetabolite or vitamin d or depigmentation or depigmenting or hydroquinone)))

2002-11-07 11:28:01

S3515

U

USPT,JPAB,EPAB,DWPI

((lisophyline or a802715 or sulphasalazine)or (alpha keto acid or pyruvic or keto octanoic or beta keto acid or acetoacetic or retinoid or retinoic or retinol or anthralin or anthranoid or dioxyanthranol or peroxide or benzoyl peroxide or minoxidil or lithium salt or antimetabolite or vitamin d or depigmentation or depigmenting or hydroquinone)) and (pharmaceutical or cosmetic)

2002-11-07 11:27:22

S3514

U

USPT,JPAB,EPAB,DWPI

(lisophyline or a802715 or sulphasalazine) or (tnf near alpha near antagonist)

2002-11-07 11:26:10

S3513

U

USPT,JPAB,EPAB,DWPI

(auranofin or skf-105809 or lactoferin) and (lisophyline or a802715 or sulphasalazine)

2002-11-07 11:02:10

S3512

U

USPT,JPAB,EPAB,DWPI

((interleukin 1 near antagonist)and (interleukin 1 near antagonist)) not ((interleukin 1 near antagonist)same (interleukin 1 near antagonist))

2002-11-07 11:00:54

S3511

U

USPT,JPAB,EPAB,DWPI

(interleukin 1 near antagonist) and (tnf near alpha near antagonist)

2002-11-07 11:00:46

S3510

U

USPT,JPAB,EPAB,DWPI

(interleukin 1 near antagonist) same (tnf near alpha near antagonist)

2002-11-07

10:58:46

S3509
U
USPT,JPAB,EPAB,DWPI
cosmetic and (((auranofin or skf-105809 or lactoferin)or (lisophyline or a802715 or
sulphasalazine)or (interleukin 1 near antagonist)or (tnf near alpha near antagonist)) and (alpha
keto acid or pyruvic or keto octanoic or beta keto acid or acetoacetic or retinoid or retinoic or
retinol or anthralin or anthranoid or dioxyanthranol or peroxide or benzoyl peroxide or minoxidil
or lithium salt or antimetabolite or vitamin d or depigmentation or depigmenting or
hydroquinone))
2002-11-07
10:22:37

S3508
U
USPT,JPAB,EPAB,DWPI
((auranofin or skf-105809 or lactoferin) or (lisophyline or a802715 or sulphasalazine) or
(interleukin 1 near antagonist) or (tnf near alpha near antagonist)) and (alpha keto acid or
pyruvic or keto octanoic or beta keto acid or acetoacetic or retinoid or retinoic or retinol or
anthralin or anthranoid or dioxyanthranol or peroxide or benzoyl peroxide or minoxidil or
lithium salt or antimetabolite or vitamin d or depigmentation or depigmenting or hydroquinone)
2002-11-07
10:22:24

S3507
U
USPT,JPAB,EPAB,DWPI
tnf near alpha near antagonist
2002-11-07
10:21:57

S3506
U
USPT,JPAB,EPAB,DWPI
interleukin 1 near antagonist
2002-11-07
10:21:23

S3505
U

USPT,JPAB,EPAB,DWPI
alpha keto acid or pyruvic or keto octanoic or beta keto acid or acetoacetic or retinoid or retinoic
or retinol or anthralin or anthranoid or dioxyanthranol or peroxide or benzoyl peroxide or
minoxidil or lithium salt or antimetabolite or vitamin d or depigmentation or depigmenting or
hydroquinone

2002-11-07 10:19:41

S3504
U
USPT,JPAB,EPAB,DWPI
lisophyline or a802715 or sulphasalazine

2002-11-07 10:15:18

S3503
U
USPT,JPAB,EPAB,DWPI
auranofin or skf-105809 or lactoferin

2002-11-07 10:14:53

S3502
U
USPT,JPAB,EPAB,DWPI
benzothiophene near carboxamide

2002-11-07 10:14:21

S3501
U
USPT,JPAB,EPAB,DWPI
furan or benzofuran

2002-11-07 10:13:52

S3500
U
USPT,JPAB,EPAB,DWPI
(loratidine or cetirizine or auranofin or lisophylline or sulphasalazine or setastine or crotamiton
or dexchlorpheniramine) or cinnarizine or cyclizine or dexchlorpheniramine or triprolidine or
phenothiazine or promethazine or alimemazine or ebastine

2002-11-07 10:13:29

S3499
U

USPT,JPAB,EPAB,DWPI

loratidine or cetirizine or auranofin or lisophylline or sulphasalazine or setastine or crotamiton or

dexchlorpheniramine

2002-11-07
10:12:31

S3498

U

USPT

alpha keto acid and beta keto acid

2002-11-07
09:59:50

```
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PASSWORD:
TERMINAL (ENTER 1, 2, 3, OR ?):2


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 NEWS  3  Apr 09   BEILSTEIN: Reload and Implementation of a New Subject Area
 NEWS  4  Apr 09   ZDB will be removed from STN
 NEWS  5  Apr 19   US Patent Applications available in IFICDB, IFIPAT, and
IFIUDB
 NEWS  6  Apr 22   Records from IP.com available in CAPLUS, HCAPLUS, and
ZCAPLUS
 NEWS  7  Apr 22   BIOSIS Gene Names now available in TOXCENTER
 NEWS  8  Apr 22   Federal Research in Progress (FEDRIP) now available
 NEWS  9  Jun 03   New e-mail delivery for search results now available
 NEWS 10  Jun 10   MEDLINE Reload
 NEWS 11  Jun 10   PCTFULL has been reloaded
 NEWS 12  Jul 02   FOREGE no longer contains STANDARDS file segment
 NEWS 13  Jul 22   USAN to be reloaded July 28, 2002;
                   saved answer sets no longer valid
 NEWS 14  Jul 29   Enhanced polymer searching in REGISTRY
 NEWS 15  Jul 30   NETFIRST to be removed from STN
 NEWS 16  Aug 08   CANCERLIT reload
 NEWS 17  Aug 08   PHARMAMarketLetter(PHARMAML) - new on STN
 NEWS 18  Aug 08   NTIS has been reloaded and enhanced
 NEWS 19  Aug 19   Aquatic Toxicity Information Retrieval (AQUIRE)
                   now available on STN
 NEWS 20  Aug 19   IFIPAT, IFICDB, and IFIUDB have been reloaded
 NEWS 21  Aug 19   The MEDLINE file segment of TOXCENTER has been reloaded
 NEWS 22  Aug 26   Sequence searching in REGISTRY enhanced
 NEWS 23  Sep 03   JAPIO has been reloaded and enhanced
 NEWS 24  Sep 16   Experimental properties added to the REGISTRY file
 NEWS 25  Sep 16   Indexing added to some pre-1967 records in CA/CAPLUS
 NEWS 26  Sep 16   CA Section Thesaurus available in CAPLUS and CA
 NEWS 27  Oct 01   CASREACT Enriched with Reactions from 1907 to 1985
 NEWS 28  Oct 21   EVENTLINE has been reloaded
 NEWS 29  Oct 24   BEILSTEIN adds new search fields
 NEWS 30  Oct 24   Nutraceuticals International (NUTRACEUT) now available on
STN
 NEWS 31  Oct 25   MEDLINE SDI run of October 8, 2002

 NEWS EXPRESS  October 14 CURRENT WINDOWS VERSION IS V6.01,
               CURRENT MACINTOSH VERSION IS V6.0a(ENG) AND V6.0Ja(JP),
               AND CURRENT DISCOVER FILE IS DATED 01 OCTOBER 2002
 NEWS HOURS    STN Operating Hours Plus Help Desk Availability
 NEWS INTER    General Internet Information
 NEWS LOGIN    Welcome Banner and News Items
 NEWS PHONE    Direct Dial and Telecommunication Network Access to STN
 NEWS WWW      CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that
specific topic.
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=> fil reg

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FULL ESTIMATED COST	0.21	0.21

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STRUCTURE FILE UPDATES: 6 NOV 2002 HIGHEST RN 471238-76-3
DICTIONARY FILE UPDATES: 6 NOV 2002 HIGHEST RN 471238-76-3

TSCA INFORMATION NOW CURRENT THROUGH MAY 20, 2002

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details:
http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf

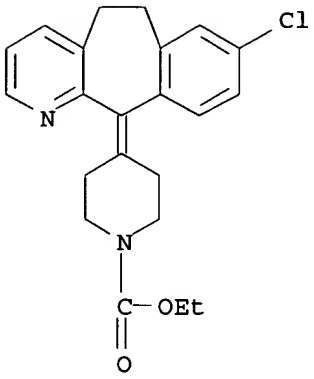
=> e loratidine/cn

E1	1	LORATADINE/CN
E2	1	LORATADINE-PSEUDOEPHEDRINE SULFATE MIXT./CN
E3	1 -->	LORATIDINE/CN
E4	1	LORATYNE/CN
E5	1	LORATYNE D/CN
E6	1	LORAX/CN
E7	1	LORAXANTHAL MONOEPOXIDE/CN
E8	1	LORAXANTHIN DIMETHYL ETHER/CN
E9	1	LORAXANTHIN MONOEPOXIDE TRIACETATE/CN
E10	1	LORAZENE/CN
E11	1	LORAZEP/CN
E12	1	LORAZEPAM/CN

=> s e3
L1 1 LORATIDINE/CN

=> d

L1 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS
RN 79794-75-5 REGISTRY
CN 1-Piperidinecarboxylic acid, 4-(8-chloro-5,6-dihydro-11H-benzo[5,6]cyclohepta[1,2-b]pyridin-11-ylidene)-, ethyl ester (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN 11H-Benzo[5,6]cyclohepta[1,2-b]pyridine, 1-piperidinecarboxylic acid deriv.
OTHER NAMES:
CN Anhissen
CN Bonalerg
CN Civeran
CN Claratyne
CN Claritin
CN Claritine
CN Clarityn
CN Clarityne
CN Cronopen
CN Fristamin
CN Histaloran
CN Klaritin
CN Lertamine
CN Lisino
CN Loracert
CN Loradex
CN Loranox
CN Lorastine
CN Loratadine
CN **Loratidine**
CN Loratyne
CN Lorfast
CN Lowadina
CN Optimin
CN Polaratyne
CN Pylor
CN Restamine
CN Sch 29851
CN Sensibit
CN Sohotin
CN Tadine
CN Velodan
CN Zeos
MF C22 H23 Cl N2 O2
CI COM
LC STN Files: ADISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CANCERLIT, CAPLUS, CASREACT, CBNB, CEN, CHEMCATS, CIN, CSCHEM, DDFU, DIOGENES, DRUGNL, DRUGPAT, DRUGU, DRUGUPDATES, EMBASE, HSDB*, IPA, MEDLINE, MRCK*, NIOSHTIC, PHAR, PHARMASEARCH, PROMT, SPECINFO, SYNTHLINE, TOXCENTER, USAN, USPAT2, USPATFULL, VETU
(*File contains numerically searchable property data)
Other Sources: WHO



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

453 REFERENCES IN FILE CA (1962 TO DATE)
12 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
459 REFERENCES IN FILE CAPLUS (1962 TO DATE)

```
=> e cetrizine/cn
E1            1     CETRIPS/CN
E2            1     CETRIZET/CN
E3            0 --> CETRIZINE/CN
E4            1     CETRORELIX/CN
E5            1     CETRORELIX ACETATE/CN
E6            1     CETROTIDE/CN
E7            1     CETUXIMAB/CN
E8            1     CETWIST TRANSCRIPTION FACTOR (CAENORHABDITIS ELEGANS GENE
HL
                    H-8)/CN
E9            1     CETYL .BETA.-AMINOCROTONATE/CN
E10           1     CETYL .GAMMA.-AMINOBUTYRATE/CN
E11           1     CETYL 1,3-DIMETHYLBUTYL ETHER/CN
E12           1     CETYL 1-NAPHTHYL ETHER/CN

=> e cetirizine/cn
E1            1     CETIPRIN NOVUM/CN
E2            1     CETIR/CN
E3            1 --> CETIRIZINE/CN
E4            1     CETIRIZINE DIHYDROCHLORIDE/CN
E5            1     CETIRIZINE HYDROCHLORIDE/CN
E6            1     CETL/CN
E7            1     CETO/CN
E8            1     CETO-2-RHODOVIBRINE/CN
E9            1     CETOBEMIDON/CN
E10           1     CETOBEMIDONE/CN
E11           1     CETOCIRE/CN
E12           1     CETOCYCLINE/CN

=> s e3
L2            1 CETIRIZINE/CN

=> d

L2   ANSWER 1 OF 1  REGISTRY  COPYRIGHT 2002 ACS
RN   83881-51-0  REGISTRY
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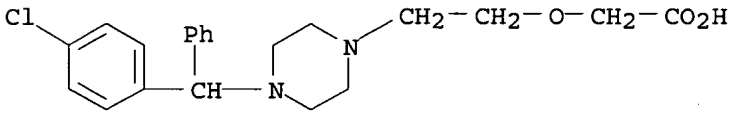
CN Acetic acid, [2-[4-[(4-chlorophenyl)phenylmethyl]-1-piperazinyl]ethoxy]-
(9CI) (CA INDEX NAME)
OTHER NAMES:
CN **Cetirizine**
FS 3D CONCORD
DR 130018-86-9
MF C21 H25 Cl N2 O3
CI COM
LC STN Files: ADISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, BIOBUSINESS,
BIOSIS, BIOTECHNO, CA, CANCERLIT, CAPLUS, CASREACT, CBNB, CEN,
CHEMCATS,
CIN, CSCHEM, DDFU, DIOGENES, DRUGNL, DRUGPAT, DRUGU, DRUGUPDATES,
EMBASE, IPA, MEDLINE, MRCK*, NIOSHTIC, PHAR, PROMT, RTECS*, SYNTHLINE,
TOXCENTER, USAN, USPAT2, USPATFULL, VETU
(*File contains numerically searchable property data)
Other Sources: WHO



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

430 REFERENCES IN FILE CA (1962 TO DATE)
9 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
435 REFERENCES IN FILE CAPLUS (1962 TO DATE)

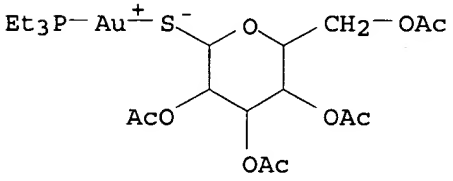
=> e auranofin/cn
E1 1 AURANAMIDE/CN
E2 1 AURANETIN/CN
E3 1 --> AURANOFIN/CN
E4 1 AURANTHINE/CN
E5 3 AURANTIA/CN
E6 1 AURANTIACIN/CN
E7 1 AURANTIACIN, DIACETATE/CN
E8 1 AURANTIACIN, DIBENZOATE/CN
E9 1 AURANTIACONE/CN
E10 1 AURANTIAMARIN/CN
E11 1 AURANTIAMIDE/CN
E12 1 AURANTIAMIDE ACETATE/CN

=> s e3
L3 1 AURANOFIN/CN

=> d

L3 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS
RN 34031-32-8 REGISTRY
CN Gold, [1-(thio-.kappa.S)-.beta.-D-glucopyranose 2,3,4,6-
tetraacetato](triethylphosphine)- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN .beta.-D-Glucopyranose, 1-thio-, 2,3,4,6-tetraacetate, gold complex
CN Gold, (1-thio-.beta.-D-glucopyranosato)(triethylphosphine)-,
2,3,4,6-tetraacetate (8CI)
CN Gold, (1-thio-.beta.-D-glucopyranose 2,3,4,6-tetraacetato-

S)(triethylphosphine)-
OTHER NAMES:
CN **Auranofin**
CN Ridaura
CN SKF 39162
CN SKF 39162D
CN [(Tetra-O-acetyl-.beta.-D-glucopyranosyl)thio](triethylphosphine)gold
MF C20 H34 Au O9 P S
CI CCS
LC STN Files: ADISNEWS, AGRICOLA, ANABSTR, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CANCERLIT, CAPLUS, CBNB, CHEMCATS, CHEMLIST, CIN, CSCHEM, DDFU, DIOGENES, DRUGPAT, DRUGU, EMBASE, GMELIN*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*, MSDS-OHS, NIOSHTIC, PHAR, PHARMASEARCH, PROMT, RTECS*, SYNTHLINE, TOXCENTER, USAN, USPATFULL, VETU
(*File contains numerically searchable property data)
Other Sources: EINECS**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)



465 REFERENCES IN FILE CA (1962 TO DATE)
9 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
466 REFERENCES IN FILE CAPLUS (1962 TO DATE)

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=> e lisophyline/cn
E1            1     LISOFYLLINE/CN
E2            1     LISOLIPIN/CN
E3            0 --> LISOPHYLINE/CN
E4            1     LISOPHYLLINE/CN
E5            1     LISORIL/CN
E6            1     LISPAMINA/CN
E7            1     LISPAMOL/CN
E8            1     LISPAMOL FUMARATE/CN
E9            1     LISPRIL/CN
E10           1     LISSAMINE AMARANTH AC/CN
E11           1     LISSAMINE BLUE/CN
E12           1     LISSAMINE BLUE 2BR/CN

=> s e4
L4            1 LISOPHYLLINE/CN

=> e
E13           1     LISSAMINE BLUE AR/CN
E14           1     LISSAMINE BLUE B/CN
E15           1     LISSAMINE BLUE BF/CN
E16           1     LISSAMINE BLUE RB/CN
E17           1     LISSAMINE FAST RED/CN
E18           1     LISSAMINE FAST RED 3G/CN
E19           1     LISSAMINE FAST RED 4G/CN
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E20          1     LISSAMINE FAST RED B/CN
E21          1     LISSAMINE FAST RED BG/CN
E22          1     LISSAMINE FAST RED BS/CN
E23          1     LISSAMINE FAST VIOLET 2B/CN
E24          1     LISSAMINE FAST YELLOW/CN

=> e l4
E1           4     L3T4.25/BI
E2           4     L3T425/BI
E3         335 --> L4/BI
E4           1     L4,1/BI
E5           2     L4.13.2/BI
E6          19     L40/BI
E7           1     L4000/BI
E8           1     L40000/BI
E9           1     L40001/BI
E10          1     L40002/BI
E11          1     L40003/BI
E12          1     L40004/BI

=> d l4

L4   ANSWER 1 OF 1  REGISTRY  COPYRIGHT 2002 ACS
RN   100324-81-0  REGISTRY
CN   1H-Purine-2,6-dione, 3,7-dihydro-1-[(5R)-5-hydroxyhexyl]-3,7-dimethyl-
     (9CI)  (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN   1H-Purine-2,6-dione, 3,7-dihydro-1-(5-hydroxyhexyl)-3,7-dimethyl-, (R)-
OTHER NAMES:
CN   CT 1501R
CN   Lisofylline
CN   Lisophylline
CN   ProTec
FS   STEREOSEARCH
MF   C13 H20 N4 O3
SR   CA
LC   STN Files:   ADISINSIGHT, ADISNEWS, AGRICOLA, BEILSTEIN*, BIOBUSINESS,
       BIOSIS, BIOTECHNO, CA, CAPLUS, CASREACT, CIN, DDFU, DRUGNL, DRUGPAT,
       DRUGU, DRUGUPDATES, EMBASE, IPA, MRCK*, PHAR, PROMT, SYNTHLINE,
       TOXCENTER, USAN, USPATFULL
         (*File contains numerically searchable property data)
     Other Sources:   WHO

Absolute stereochemistry.  Rotation (-).
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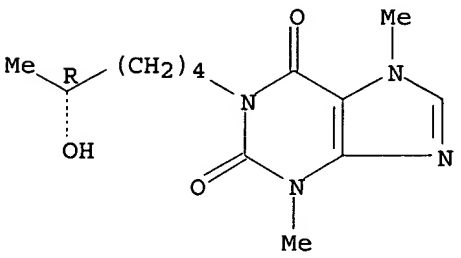



```
**PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT**
```

63 REFERENCES IN FILE CA (1962 TO DATE)
63 REFERENCES IN FILE CAPLUS (1962 TO DATE)

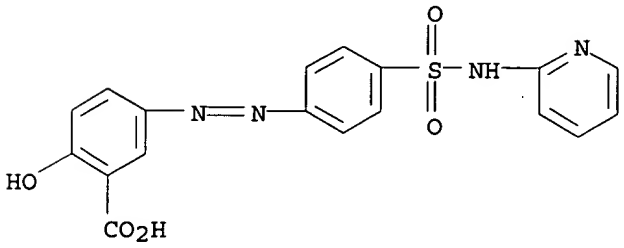
```
=> e sulphasalazine/cn
E1            1     SULPHARSENOLUM/CN
E2            1     SULPHARSIDE/CN
E3            1 --> SULPHASALAZINE/CN
E4            1     SULPHASIL/CN
E5            1     SULPHASOLUCIN/CN
E6            1     SULPHASOMIDINE/CN
E7            1     SULPHASOMIZOLE/CN
E8            1     SULPHAST GREEN/CN
E9            1     SULPHATASE (SALMONELLA ENTERICA TYPHI STRAIN CT18 GENE
STY24
                    66)/CN
E10           1     SULPHATE/CN
E11           1     SULPHATE ABC TRANSPORTER (NEISSERIA MENINGITIDIS STRAIN
C311
                    + CLONE NMB0580 GENE PHO2-91 HOMOLOG)/CN
E12           1     SULPHATE ABC TRANSPORTER (NEISSERIA MENINGITIDIS STRAIN
C311
                    + CLONE NMB0880 GENE PHO2-90 HOMOLOG)/CN

=> s e3
L5            1 SULPHASALAZINE/CN

=> d

L5   ANSWER 1 OF 1  REGISTRY  COPYRIGHT 2002 ACS
RN   599-79-1  REGISTRY
CN   Benzoic acid, 2-hydroxy-5-[[4-[(2-pyridinylamino)sulfonyl]phenyl]azo]-
     (9CI)  (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN   Salicylic acid, 5-[[p-(2-pyridylsulfamoyl)phenyl]azo]- (6CI, 7CI, 8CI)
OTHER NAMES:
CN   2-Hydroxy-5-((4-((2-pyridinylamino)sulfonyl)phenyl)azo)benzoic acid
CN   5-[4-(2-Pyridylsulfamyl)phenylazo]-2-hydroxybenzoic acid
CN   5-[p-(2-Pyridylsulfamoyl)phenylazo]salicylic acid
CN   5-[p-(2-Pyridylsulfamyl)phenylazo]salicylic acid
CN   Azopyrin
CN   Azopyrine
CN   Azulfidine
CN   Benzosulfa
CN   Reupirin
CN   Salazopyridin
CN   Salazopyrin
CN   Salazopyrine
CN   Salazosulfapyridin
CN   Salazosulfapyridine
CN   Salicylazosulfapyridine
CN   Salisulf
CN   Sulfasalazin
CN   Sulfasalazine
CN   Sulphasalazine
FS   3D CONCORD
MF   C18 H14 N4 O5 S
CI   COM
LC   STN Files:   ADISNEWS, AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS,
BIOSIS,
```

BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB, CEN, CHEMCATS,
CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DIOGENES, DRUGNL, DRUGU,
DRUGUPDATES,
EMBASE, HODOC*, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*,
NIOSHTIC, PHAR, PHARMASEARCH, PIRA, PROMT, RTECS*, SPECINFO, TOXCENTER,
USAN, USPATFULL, VETU
(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

899 REFERENCES IN FILE CA (1962 TO DATE)
32 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
903 REFERENCES IN FILE CAPLUS (1962 TO DATE)
12 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

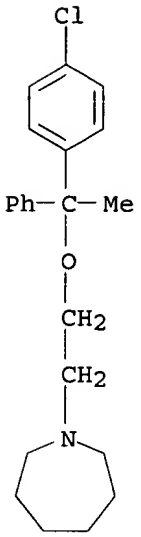
```
=> e setastine/cn
E1            1     SETARON GOLDEN YELLOW RL/CN
E2            1     SETARON RED BL/CN
E3            1 --> SETASTINE/CN
E4            1     SETAVIN CAW/CN
E5            1     SETAVIN KS/CN
E6            1     SETAVIN PE/CN
E7            1     SETAVONE C/CN
E8            1     SETAVONE O/CN
E9            1     SETAVONE ST/CN
E10           1     SETAZINDOL/CN
E11           1     SETAZINE/CN
E12           1     SETCREASIN/CN

=> s e3
L6            1 SETASTINE/CN

=> d

L6   ANSWER 1 OF 1  REGISTRY  COPYRIGHT 2002 ACS
RN   64294-95-7  REGISTRY
CN   1H-Azepine, 1-[2-[1-(4-chlorophenyl)-1-phenylethoxy]ethyl]hexahydro-
(9CI)
     (CA INDEX NAME)
OTHER NAMES:
CN   Setastine
FS   3D CONCORD
MF   C22 H28 Cl N O
CI   COM
```

LC STN Files: BEILSTEIN*, BIOSIS, CA, CAPLUS, CIN, DDFU, DRUGU, EMBASE,
MEDLINE, MRCK*, SYNTHLINE, TOXCENTER, USAN, USPATFULL
(*File contains numerically searchable property data)
Other Sources: WHO



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

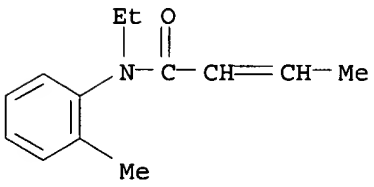
25 REFERENCES IN FILE CA (1962 TO DATE)
3 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
25 REFERENCES IN FILE CAPLUS (1962 TO DATE)

```
=> e crotamiton/cn
E1            1     CROTAMINE, PRO- (CROTALUS DURISSUS TERRIFICUS CLONE PCM31B
R
                    EDUCED)/CN
E2            1     CROTAMINE, PRO- (CROTALUS DURISSUS TERRIFICUS CLONE PCM45B
R
                    EDUCED)/CN
E3            1 --> CROTAMITON/CN
E4            1     CROTAMITON-DELTAMETHRIN MIXT./CN
E5            1     CROTAMITON-PERMETHRIN MIXT./CN
E6            1     CROTAMITON-RESMETHRIN MIXT./CN
E7            1     CROTAMITON-TETRAMETHRIN MIXT./CN
E8            1     CROTAMITONE/CN
E9            1     CROTAN BK/CN
E10           1     CROTANANIC ACID/CN
E11           1     CROTANANIC ACID LACTONE/CN
E12           1     CROTANANINE/CN

=> s e3
L7            1 CROTAMITON/CN

=> d
```

L7 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS
RN 483-63-6 REGISTRY
CN 2-Butenamide, N-ethyl-N-(2-methylphenyl)- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN o-Crotonotoluidide, N-ethyl- (6CI, 7CI, 8CI)
OTHER NAMES:
CN Crotalgin
CN **Crotamiton**
CN Crotamitone
CN Crotonyl N-ethyl-o-toluidine
CN Eurax
CN Euraxil
CN N-Crotonyl-N-ethyl-o-toluidine
CN N-Ethyl-o-crotonotoluidide
CN Veteusan
FS 3D CONCORD
MF C13 H17 N O
CI COM
LC STN Files: ADISNEWS, AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS,
BIOTECHNO, CA, CABA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB, CEN, CHEMCATS, CHEMLIST, CIN, CSCHEM, DDFU, DIOGENES, DRUGU, EMBASE, IPA, MEDLINE, MRCK*, PHARMASEARCH, PROMT, RTECS*, SPECINFO, TOXCENTER, USAN, USPATFULL, VETU
(*File contains numerically searchable property data)
Other Sources: EINECS**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

215 REFERENCES IN FILE CA (1962 TO DATE)
3 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
215 REFERENCES IN FILE CAPLUS (1962 TO DATE)
2 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

=> e dexchlorpheniramine/cn
E1 1 DEXCEL CLEAR CONC S 380/CN
E2 1 DEXCHLOROPHENIRAMINE MALEATE/CN
E3 1 --> DEXCHLORPHENIRAMINE/CN
E4 1 DEXCHLORPHENIRAMINE HYDROGEN MALEATE/CN
E5 1 DEXCHLORPHENIRAMINE MALEATE/CN
E6 1 DEXCLAMOL/CN
E7 1 DEXCLAMOL HYDROBROMIDE/CN
E8 1 DEXCLAMOL HYDROCHLORIDE/CN
E9 1 DEXCO 2518/CN
E10 1 DEXCO 4111/CN

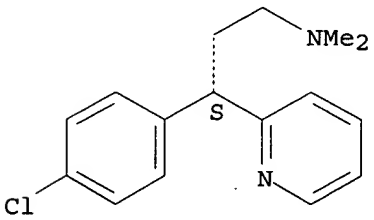
```
E11           1     DEXCO 506A/CN
E12           1     DEXCO 8508D/CN

=> s e3
L8            1 DEXCHLORPHENIRAMINE/CN

=> d

L8   ANSWER 1 OF 1  REGISTRY  COPYRIGHT 2002 ACS
RN   25523-97-1  REGISTRY
CN   2-Pyridinepropanamine, .gamma.-(4-chlorophenyl)-N,N-dimethyl-,
(.gamma.S)-
     (9CI)  (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN   2-Pyridinepropanamine, .gamma.-(4-chlorophenyl)-N,N-dimethyl-, (S)-
CN   Pyridine, 2-[p-chloro-.alpha.-[2-(dimethylamino)ethyl]benzyl]-, (S)-(+)-
     (8CI)
OTHER NAMES:
CN   (+)-Chlorpheniramine
CN   d-Chlorpheniramine
CN   Dexchlorpheniramine
CN   S-(+)-Chlorpheniramine
CN   S-Chlorpheniramine
FS   STEREOSEARCH
DR   12167-38-3, 301-24-6
MF   C16 H19 Cl N2
CI   COM
LC   STN Files:   ADISNEWS, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS,
       BIOTECHNO, CA, CANCERLIT, CAPLUS, CASREACT, CBNB, CHEMLIST, CIN, DDFU,
       DRUGU, EMBASE, HSDB*, IPA, MEDLINE, MRCK*, PHAR, PROMT, SPECINFO,
       TOXCENTER, USAN, USPAT2, USPATFULL, VETU
         (*File contains numerically searchable property data)
     Other Sources:   EINECS**, WHO
         (**Enter CHEMLIST File for up-to-date regulatory information)

Absolute stereochemistry.  Rotation (+).
```

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**PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT**

             217 REFERENCES IN FILE CA (1962 TO DATE)
               4 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
             218 REFERENCES IN FILE CAPLUS (1962 TO DATE)


=> e benzyloxy(l)methoxy(l)tetrazol(l)benzothiophene(l)carboxamide
E1            1     BENZYLOXPHENYLACETIC/BI
E2        23201     BENZYLOXY/BI
```

```
E3            0 -->
BENZYLOXY(L)METHOXY(L)TETRAZOL(L)BENZOTHIOPHENE(L)CARBOXAMID
                    E/BI
E4           43     BENZYLOXYACET/BI
E5            2     BENZYLOXYACETALDEHYDE/BI
E6            6     BENZYLOXYACETAMIDE/BI
E7            2     BENZYLOXYACETANILIDE/BI
E8            5     BENZYLOXYACETATE/BI
E9           10     BENZYLOXYACETIC/BI
E10          26     BENZYLOXYACETO/BI
E11           2     BENZYLOXYACETOACETIC/BI
E12           1     BENZYLOXYACETONE/BI

=> e benzyloxy(l)methoxy(l)tetrazol(l)benzothiophene(l)carboxamide/cn
E1            1     BENZYLOXY TERT-BUTYL NITROXIDE/CN
E2            1     BENZYLOXY(ETHYL)AMINE/CN
E3            0 -->
BENZYLOXY(L)METHOXY(L)TETRAZOL(L)BENZOTHIOPHENE(L)CARBOXAMID
                    E/CN
E4            1     BENZYLOXY(PHENYLSULFONYL)METHANE/CN
E5            1     BENZYLOXY(TERT-BUTYL)DIMETHYLSILANE/CN
E6            1     BENZYLOXY, .ALPHA.,.ALPHA.-DIMETHYL-/CN
E7            1     BENZYLOXY, .ALPHA.-ETHYL-.ALPHA.-METHYL-/CN
E8            1     BENZYLOXY, .ALPHA.-METHYL-/CN
E9            1     BENZYLOXY, DIHYDROXY-/CN
E10           1     BENZYLOXY, P-METHYL-/CN
E11           1     BENZYLOXY-1-BROMO-2-FLUOROBENZENE/CN
E12           1     BENZYLOXY-1-NAPHTHYLPHENYLSILANE/CN

=> log y
COST IN U.S. DOLLARS                          SINCE FILE      TOTAL
                                                 ENTRY    SESSION
FULL ESTIMATED COST                              47.30      47.51

STN INTERNATIONAL LOGOFF AT 09:35:44 ON 07 NOV 2002
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